

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP092619\
 Data File : BP000445.D
 Acq On : 27 Sep 2019 03:26
 Operator : JU
 Sample : K4663-09
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 OK-02-092419

Quant Time: Sep 27 06:34:18 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP091619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 19 14:44:51 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	245194	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	890012	20.00	ng	0.00
39) Acenaphthene-d10	9.83	164	487213	20.00	ng	0.00
64) Phenanthrene-d10	11.33	188	874611	20.00	ng	0.00
76) Chrysene-d12	14.00	240	742713	20.00	ng	0.00
87) Perylene-d12	15.49	264	810090	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.43	112	1861634	131.07	ng	0.02
7) Phenol-d6	6.43	99	2355517	135.30	ng	0.01
23) Nitrobenzene-d5	7.35	82	1359816	98.58	ng	0.00
42) 2,4,6-Tribromophenol	10.63	330	741471	153.43	ng	0.00
45) 2-Fluorobiphenyl	9.15	172	2792032	103.14	ng	0.00
79) Terphenyl-d14	12.94	244	3391616	95.80	ng	0.00
Target Compounds						
10) Phenol	6.45	94	144051	7.283	ng	97
50) Dimethylphthalate	9.55	163	389914	11.746	ng	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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